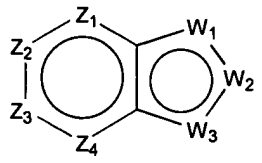
AMENDMENTS TO THE CLAIMS

Please amend the claims as follows:

1. (CURRENTLY AMENDED) A compound of formula (I):



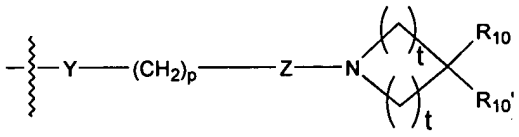
(I)

wherein:

Z_1 is CR_1 ~~or N~~, Z_2 is CR_2 ~~or N~~, Z_3 is CR_3 ~~or N~~, and Z_4 is CR_4 ~~or N, where no more than two of Z_1, Z_2, Z_3 and Z_4 are N~~;

W_1 is O, S, or NR_5, ~~one of~~ W_2 ~~and W_3~~ is N or CR_6, and ~~the other of W_2 and~~ W_3 is CG; W_1 is NG, W_2 is CR_5 or N, and W_3 is CR_6 or N; ~~or W_1 and W_3 are N, and W_2 is NG~~;

G is of formula (II):



(II)

Y is O, S, CHOH, -NHC(O)-, -C(O)NH-, -C(O)-, -OC(O)-, -(O)CO-, $-NR_7$-, -CH=N-, or absent;

p is 1, 2, 3, 4 or 5;

Z is CR_8R_9 or absent;

each t is 1, 2, or 3;

each R_1, R_2, R_3, and R_4, independently, is H, amino, hydroxyl, halo, or straight- or branched-chain C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, C_{1-6} heteroalkyl, C_{1-6} haloalkyl, -CN, $-CF_3$, $-OR_{11}$, $-COR_{11}$, $-NO_2$, $-SR_{11}$, $-NHC(O)R_{11}$, $-C(O)NR_{12}R_{13}$, $-NR_{12}R_{13}$, $-NR_{11}C(O)NR_{12}R_{13}$, $-SO_2NR_{12}R_{13}$, $-OC(O)R_{11}$, $-O(CH_2)_qNR_{12}R_{13}$, or $-(CH_2)_qNR_{12}R_{13}$, where q is an integer from 2 to 6, or R_1 and R_2 together form -NH-N=N- or R_3 and R_4 together form -NH-N=N-;

each R_5, R_6, and R_7, independently, is H, C_{1-6} alkyl; formyl; C_{3-6} cycloalkyl; C_{5-6} aryl, optionally substituted with halo or C_{1-6} alkyl; or C_{5-6} heteroaryl, optionally substituted with halo or C_{1-6} alkyl;

each R_8 and R_9, independently, is H or straight- or branched-chain C_{1-8} alkyl;

R_{10} is H, straight- or branched-chain C_{1-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, C_{1-8} alkylidene, C_{1-8} alkoxy, C_{1-8} heteroalkyl, C_{1-8} aminoalkyl, C_{1-8} haloalkyl, C_{1-8} alkoxycarbonyl, C_{1-8} hydroxyalkoxy, C_{1-8} hydroxyalkyl, -SH, C_{1-8} alkylthio, -O-CH_2-C_{5-6} aryl, -C(O)-C_{5-6} aryl substituted with C_{1-3} alkyl or halo, C_{5-6} aryl, C_{5-6} cycloalkyl, C_{5-6} heteroaryl, C_{5-6} heterocycloalkyl, -$NR_{12}R_{13}$, -$C(O)NR_{12}R_{13}$, -$NR_{11}C(O)NR_{12}R_{13}$, -$CR_{11}R_{12}R_{13}$, -$OC(O)R_{11}$, -$(O)(CH_2)_sNR_{12}R_{13}$ or -$(CH_2)_sNR_{12}R_{13}$, s being an integer from 2 to 8;

R_{10}' is H, straight- or branched-chain C_{1-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, C_{1-8} alkylidene, C_{1-8} alkoxy, C_{1-8} heteroalkyl, C_{1-8} aminoalkyl, C_{1-8} haloalkyl, C_{1-8} alkoxycarbonyl, C_{1-8} hydroxyalkoxy, C_{1-8} hydroxyalkyl, or C_{1-8} alkylthio;

each R_{11}, independently, is H, straight- or branched-chain C_{1-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, C_{2-8} heteroalkyl, C_{2-8} aminoalkyl, C_{2-8} haloalkyl, C_{1-8} alkoxycarbonyl, C_{2-8} hydroxyalkyl, -C(O)-C_{5-6} aryl substituted with C_{1-3} alkyl or halo, C_{5-6} aryl, C_{5-6} heteroaryl, C_{5-6} cycloalkyl, C_{5-6} heterocycloalkyl, -$C(O)NR_{12}R_{13}$, -$CR_5R_{12}R_{13}$, -$(CH_2)_tNR_{12}R_{13}$, t is an integer from 2 to 8; and

each R_{12} and R_{13}, independently, is H, C_{1-6} alkyl; C_{3-6} cycloalkyl; C_{5-6} aryl, optionally substituted with halo or C_{1-6} alkyl; or C_{5-6} heteroaryl, optionally substituted with halo or C_{1-6} alkyl; or R_{12} and R_{13} together form a cyclic structure;

or a pharmaceutically acceptable salt, ester or prodrug thereof.

2. (ORIGINAL) The compound of claim 1, wherein each t is 2 and R_{10} is straight- or branched-chain C_{2-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, C_{1-8} alkylidene, C_{1-8} alkoxy, or C_{1-8} heteroalkyl.

3. (ORIGINAL) The compound of claim 2, wherein R_{10} is n-butyl.

4. (CANCELED)

5. (CURRENTLY AMENDED) The compound of claim 42, wherein each R_1, R_2, R_3, and R_4, independently, is H, hydroxyl, halo, C_{1-6}heteroalkyl, CF_3, -NO_2, or

straight- or branched-chain C_{1-6} alkyl, or R_1 and R_2 together form -NH-N=N- or R_3 and R_4 together form -NH-N=N-.

6. (ORIGINAL) The compound of claim 2, wherein Y is absent or O, p is 0, 1, 2 or 3, and R_8 and R_9 are H.

7. (ORIGINAL) The compound of claim 6, wherein Z is absent, Y is absent and p is 3.

8. (ORIGINAL) The compound of claim 7, wherein R_{10} is n-butyl.

9-16. (CANCELED)

17. (CURRENTLY AMENDED) The compound of claim 1, wherein the compound is:

~~2-(3-(4-*n*-butylpiperidine-1-yl)-propyl)-benzothiazole;~~

~~2-(3-(4-*n*-butylpiperidine-1-yl)-propyl)-benzooxazole;~~

~~4,5-difluoro-2-(3-(4-*n*-butylpiperidine-1-yl)-propyl)-1*H*-benzoimidazole;~~

~~6-fluoro-5-nitro-2-(3-(4-*n*-butylpiperidine-1-yl)-propyl)-1*H*-benzoimidazole;~~

~~5-tert-butyl-2-(3-(4-*n*-butylpiperidine-1-yl)-propyl)-1*H*-benzoimidazole;~~

~~5-chloro-6-methyl-2-(3-(4-*n*-butylpiperidine-1-yl)-propyl)-1*H*-benzoimidazole;~~

~~4,6-difluoro-2-(3-(4-*n*-butylpiperidine-1-yl)-propyl)-1*H*-benzoimidazole;~~

~~2-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-imidazo[*4,5-c*]pyridine;~~

~~8-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-9*H*-purine;~~

~~7-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-3,8-dihydro-imidazo[*4',5':3,4*]benzo[*1,2-d*][*1,2,3*]triazole;~~

~~2-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-3a,4,5,6,7,7a-hexahydro-1*H*-benzoimidazole;~~

1-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indole;

1-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-benzoimidazole;

3-methyl-1-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indole;

5-bromo-1-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indole;

3-formyl-1-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indole;

7-bromo-1-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indole;

~~1-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indazole;~~

3-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-benzo[*d*]isoxazole;

3-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indole;

~~4-nitro-2-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-benzoimidazole;~~

~~5-nitro-2-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-benzoimidazole;~~

~~4-hydroxy-2-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-benzoimidazole;~~

~~2-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-benzoimidazole;~~

~~4-methyl-2-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-benzoimidazole;~~

3-(2-(4-*n*-butylpiperidine)-1-yl-ethyl)-1*H*-indole;

3-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indazole;

3-(2-(4-n-butylpiperidine)-ethoxy)-7-methyl-benzo[*d*]isoxazole;

1-(3-(4-~~M~~methylpiperidine)-1-yl-propyl)-1H-indazole;

1-(3-(4-~~P~~pentylpiperidine)-1-yl-propyl)-1H-indazole;

1-(3-(4-~~P~~propylpiperidine)-1-yl-propyl)-1H-indazole;

1-(3-(4-(3-~~M~~methyl-butyl)-piperidine)-1-yl-propyl)-1H-indazole

1-(3-(4-~~P~~pentylidene-piperidine)-1-yl-propyl)-1H-indazole;

1-(3-(4-~~P~~propylidene-piperidine)-1-yl-propyl)-1H-indazole

1-~~B~~benzo[*b*]thiophen-2-yl-4-(4-butylpiperidin-1-yl)-butan-1-one

4-(4-~~B~~butylpiperidin-1-yl)-1-(3-methyl-benzofuran-2-yl)-butan-1-one;

4-(4-~~B~~butylpiperidin-1-yl)-1-(5-fluoro-3-methyl-benzo[*b*]thiophen-2-yl)-butan-1-one;

1-~~B~~benzofuran-2-yl-4-(4-butylpiperidin-1-yl)-butan-1-one;

1-(3-~~B~~bromo-benzo[*b*]thiophen-2-yl)-4-(4-butylpiperidin-1-yl)-butan-1-one

1-(3-~~B~~benzo[*b*]thiophen-2-yl-propyl)-4-butylpiperidine;

1-(3-~~B~~benzofuran-2-yl-propyl)-4-butylpiperidine;

4-~~B~~butyl-1-[3-(3-methyl-benzofuran-2-yl)-propyl]-piperidine;

4-~~B~~butyl-1-[3-(5-fluoro-3-methyl-benzo[*b*]thiophen-2-yl)-propyl]-piperidine;

2-(3-~~I~~iodo-propyl)-benzo[*b*]thiophene;

1-(3-~~B~~benzo[*b*]thiophen-2-yl-propyl)-4-methylpiperidine

1-(3-~~B~~benzo[*b*]thiophen-2-yl-propyl)-4-benzylpiperidine;

1-(3-~~B~~benzo[*b*]thiophen-2-yl-propyl)-4-(2-methoxy-phenyl)-piperidine;

~~2-(3-Bromopropyl)-2H-benzotriazole;~~

~~2-[3-(4-Butylpiperidin-1-yl)-propyl]-2H-benzotriazole;~~

~~1-(3-Bromopropyl)-1H-benzotriazole;~~

1-[3-(4-~~B~~butylpiperidin-1-yl)-propyl]-1H-benzotriazole;

1-[3-(4-~~B~~butylpiperidin-1-yl)-propyl]-1*H*-indole-3-carbaldehyde;

{1-[3-(4-~~B~~butylpiperidin-1-yl)-propyl]-1H-indol-3-yl}-methanol;

1-[3-(4-~~B~~butylpiperidin-1-yl)-propyl]-2-phenyl-1*H*-benzoimidazole;

1-[3-(4-~~B~~butylpiperidin-1-yl)-propyl]-3-chloro-1*H*-indazole;

1-[3-(4-~~B~~butylpiperidin-1-yl)-propyl]-6-nitro-1*H*-indazole;

~~Benzo[*d*]isoxazol-3-ol;~~

~~3-(2-Chloroethoxy)-benzo[*d*]isoxazole;~~

3-[2-(4-~~B~~butylpiperidin-1-yl)-ethoxy]-benzo[*d*]isoxazol;

~~3-(1H-I~~indol-3-yl)-propan-1-ol;

3-[3-(4-~~B~~butyl-piperidin-1-yl)-propyl]- 1*H*-indole hydrochloride;

~~4-(4-Butylpiperidine-1-yl)-butyric acid methyl ester;~~

~~2-[3-(4-Butylpiperidin-1-yl)-propyl]-1-methyl-1H-benzimidazole;~~

1H-~~I~~indazole-3-carboxylic acid (2-(4-butylpiperidin)-1-yl-ethyl)-amide;

1-[3-(4-~~B~~butylpiperidin-1-yl)-propyl]-5-nitro-1H-indazole;

~~2-[3-(4-butylpiperidin-1-yl)-propyl]-5-nitro-2H-indazole;~~

1-[3-(4-~~B~~butyl-piperidin-1-yl)-propyl]-2-methyl-1*H*-indole;

1-{1-[3-(4-~~B~~butyl-piperidin-1-yl)-propyl]-1H-indol-3-yl}-ethanone;

{1-[3-(4-~~B~~butyl-piperidin-1-yl)-propyl]-1H-indol-3-yl}-acetonitrile;

1-[3-(4-~~B~~butyl-piperidin-1-yl)-propyl]-1*H*-indole -3-carbonitrile;

1-[3-(4-~~B~~butyl-piperidin-1-yl)-propyl]-5,6-dimethyl-1*H*-benzoimidazole;

1-[3-(4-~~B~~butyl-piperidin-1-yl)-propyl]-5(6)-dimethyl-1*H*-benzoimidazole;

1-[3-(4-~~B~~butyl-piperidin-1-yl)-propyl]-5-methoxy-1*H*-benzoimidazole;

{1-[3-(4-~~B~~butyl-piperidin-1-yl)-propyl]-1H-benzoimidazol-2-yl}-methanol;

1-[3-(4-~~B~~butyl-piperidin-1-yl)-propyl]-2-trifuoromethyl-1*H*-benzoimidazole;

~~(2-Trimethylstannanyl-phenyl)-carbamic acid *tert*-butyl ester;~~

~~[2-(4-Chloro-butyryl)-phenyl]-carbamic acid *tert*-butyl ester;~~

~~{2-[4-(4-Butyl-piperidine-1-yl)-butyryl]-phenyl}-carbamic acid *tert*-butyl ester;~~

3-[3-(4-~~B~~butyl-piperidine-1-yl)-propyl]-1H-indazole, HCl;

3-[3-(4-~~B~~butyl-piperidine-1-yl)-propyl]-5-nitro-1H-indazole;

3-[3-(4-~~B~~butyl-piperidine-1-yl)-propyl]-5,7-dinitro-1H-indazole;

~~4-(4-Butyl-piperidin-1-yl)-1-(2-metylsulfanyl-phenyl)-butan-1-one;~~

3-[3-(4-~~B~~butyl-piperidin-1-yl)-propyl]-benzo[*d*]isothiazole;

3-[3-(4-~~B~~butyl-piperidin-1-yl)-propyl]-5-methoxy-1H-indazole;

3-[3-(4-~~B~~butyl-piperidin-1-yl)-propyl]-4-methoxy-1H-indazole

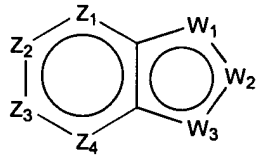
3-[3-(4-~~B~~butyl-piperidin-1-yl)-propyl]-6-methoxy-1H-indazole;

3-[3-(4-~~B~~butyl-piperidin-1-yl)-propyl]-1H-indazole-4-ol ~~(53MF51)~~;

3-[3-(4-~~B~~butyl-piperidin-1-yl)-propyl]-1H-indazole-6-ol ~~(53MF52)~~; or

3-[3-(4-~~B~~butyl-piperidin-1-yl)-propyl]-1H-indazole-5-ol.

18. (CURRENTLY AMENDED) A pharmaceutical composition comprising an effective amount of a compound of formula (I):



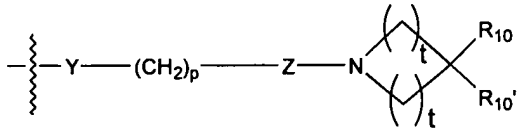
(I)

wherein:

Z_1 is CR_1 ~~or N~~, Z_2 is CR_2 ~~or N~~, Z_3 is CR_3 ~~or N~~, and Z_4 is CR_4 ~~or N, where no more than two of Z_1, Z_2, Z_3 and Z_4 are N~~;

W_1 is O, S, or NR_5, ~~one of~~ W_2 ~~and W_3~~ is N or CR_6, and ~~the other of W_2 and~~ W_3 is CG; W_1 is NG, W_2 is CR_5 or N, and W_3 is CR_6 or N; ~~or W_1 and W_3 are N, and W_2 is NG~~;

G is of formula (II):



(II)

Y is O, S, CHOH, -NHC(O)-, -C(O)NH-, -C(O)-, -OC(O)-, -(O)CO-, $-NR_7$-, -CH=N-, or absent;

p is 1, 2, 3, 4 or 5;

Z is CR_8R_9 or absent;

each t is 1, 2, or 3;

each R_1, R_2, R_3, and R_4, independently, is H, amino, hydroxyl, halo, or straight- or branched-chain C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, C_{1-6} heteroalkyl, C_{1-6} haloalkyl, -CN, -CF_3, -OR_{11}, -COR_{11}, -NO_2, -SR_{11}, -$NHC(O)R_{11}$, -$C(O)NR_{12}R_{13}$, -$NR_{12}R_{13}$, -$NR_{11}C(O)NR_{12}R_{13}$, -$SO_2NR_{12}R_{13}$, -$OC(O)R_{11}$, -$O(CH_2)_qNR_{12}R_{13}$, or -$(CH_2)_qNR_{12}R_{13}$, where q is an integer from 2 to 6, or R_1 and R_2 together form -NH-N=N- or R_3 and R_4 together form -NH-N=N-;

each R_5, R_6, and R_7, independently, is H, C_{1-6} alkyl; formyl; C_{3-6} cycloalkyl; C_{5-6} aryl, optionally substituted with halo or C_{1-6} alkyl; or C_{5-6} heteroaryl, optionally substituted with halo or C_{1-6} alkyl;

each R_8 and R_9, independently, is H or straight- or branched-chain C_{1-8} alkyl;

R_{10} is H, straight- or branched-chain C_{1-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, C_{1-8} alkylidene, C_{1-8} alkoxy, C_{1-8} heteroalkyl, C_{1-8} aminoalkyl, C_{1-8} haloalkyl, C_{1-8} alkoxycarbonyl, C_{1-8} hydroxyalkoxy, C_{1-8} hydroxyalkyl, -SH, C_{1-8} alkylthio, -O-CH_2-C_{5-6} aryl, -C(O)-C_{5-6} aryl substituted with C_{1-3} alkyl or halo, C_{5-6} aryl, C_{5-6} cycloalkyl, C_{5-6} heteroaryl, C_{5-6} heterocycloalkyl, -$NR_{12}R_{13}$, -$C(O)NR_{12}R_{13}$, -$NR_{11}C(O)NR_{12}R_{13}$, -$CR_{11}R_{12}R_{13}$, -$OC(O)R_{11}$, -$(O)(CH_2)_SNR_{12}R_{13}$ or -$(CH_2)_SNR_{12}R_{13}$, s being an integer from 2 to 8;

R_{10}' is H, straight- or branched-chain C_{1-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, C_{1-8} alkylidene, C_{1-8} alkoxy, C_{1-8} heteroalkyl, C_{1-8} aminoalkyl, C_{1-8} haloalkyl, C_{1-8} alkoxycarbonyl, C_{1-8} hydroxyalkoxy, C_{1-8} hydroxyalkyl, or C_{1-8} alkylthio;

each R_{11}, independently, is H, straight- or branched-chain C_{1-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, C_{2-8} heteroalkyl, C_{2-8} aminoalkyl, C_{2-8} haloalkyl, C_{1-8} alkoxycarbonyl, C_{2-8} hydroxyalkyl, -C(O)-C_{5-6} aryl substituted with C_{1-3} alkyl or halo, C_{5-6} aryl, C_{5-6} heteroaryl, C_{5-6} cycloalkyl, C_{5-6} heterocycloalkyl, -$C(O)NR_{12}R_{13}$, -$CR_5R_{12}R_{13}$, -$(CH_2)_tNR_{12}R_{13}$, t is an integer from 2 to 8; and

each R_{12} and R_{13}, independently, is H, C_{1-6} alkyl; C_{3-6} cycloalkyl; C_{5-6} aryl, optionally substituted with halo or C_{1-6} alkyl; or C_{5-6} heteroaryl, optionally substituted with halo or C_{1-6} alkyl; or R_{12} and R_{13} together form a cyclic structure;

or a pharmaceutically acceptable salt, ester or prodrug thereof.

19. (ORIGINAL) A pharmaceutical composition of Claim 18, wherein each t is 2 and R_{10} is straight- or branched-chain C_{2-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, C_{1-8} alkylidene, C_{1-8} alkoxy, or C_{1-8} heteroalkyl.

20. (ORIGINAL) A pharmaceutical composition of Claim 19, wherein R_{10} is n-butyl.

21. (CANCELED)

22. (CURRENTLY AMENDED) A pharmaceutical composition of Claim ~~21~~19, wherein each R_1, R_2, R_3, and R_4, independently, is H, hydroxyl, halo, C_{1-6}heteroalkyl, CF_3, $-NO_2$, or straight- or branched-chain C_{1-6} alkyl, or R_1 and R_2 together form -NH-N=N- or R_3 and R_4 together form -NH-N=N-.

23. (ORIGINAL) A pharmaceutical composition of Claim 19, wherein Y is absent or O, p is 0, 1, 2 or 3, and R_8 and R_9 are H.

24. (ORIGINAL) A pharmaceutical composition of Claim 23, wherein Z is absent, Y is absent and p is 3.

25. (ORIGINAL) A pharmaceutical composition of Claim 24, wherein R_{10} is n-butyl.

26-33. (CANCELED)

34. (CURRENTLY AMENDED) A pharmaceutical composition of Claim 19, wherein the compound is:

~~2-(3-(4-*n*-butylpiperidine-1-yl)-propyl)-benzothiazole;~~

~~2-(3-(4-*n*-butylpiperidine-1-yl)-propyl)-benzooxazole;~~

~~4,5-difluoro-2-(3-(4-*n*-butylpiperidine-1-yl)-propyl)-1*H*-benzoimidazole;~~

~~6-fluoro-5-nitro-2-(3-(4-*n*-butylpiperidine-1-yl)-propyl)-1*H*-benzoimidazole;~~

~~5-tert-butyl-2-(3-(4-*n*-butylpiperidine-1-yl)-propyl)-1*H*-benzoimidazole;~~

~~5-chloro-6-methyl-2-(3-(4-*n*-butylpiperidine-1-yl)-propyl)-1*H*-benzoimidazole;~~

~~4,6-difluoro-2-(3-(4-*n*-butylpiperidine-1-yl)-propyl)-1*H*-benzoimidazole;~~

~~2-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-imidazo[*4,5-c*]pyridine;~~

~~8-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-9*H*-purine;~~

~~7-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-3,8-dihydro-imidazo[*4',5':3,4*]benzo[*1,2-d*][*1,2,3*]triazole;~~

~~2-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-3a,4,5,6,7,7a-hexahydro-1*H*-benzoimidazole;~~

1-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indole;

1-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-benzoimidazole;

3-methyl-1-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indole;

5-bromo-1-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indole;

3-formyl-1-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indole;

7-bromo-1-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indole;

~~1-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indazole;~~

3-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-benzo[*d*]isoxazole;

3-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indole;

~~4-nitro-2-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-benzoimidazole;~~

~~5-nitro-2-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-benzoimidazole;~~

~~4-hydroxy-2-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-benzoimidazole;~~

~~2-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-benzoimidazole;~~

~~4-methyl-2-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-benzoimidazole;~~

3-(2-(4-*n*-butylpiperidine)-1-yl-ethyl)-1*H*-indole;

3-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indazole;

3-(2-(4-n-butylpiperidine)-ethoxy)-7-methyl-benzo[*d*]isoxazole;

1-(3-(4-methylpiperidine)-1-yl-propyl)-1H-indazole;

1-(3-(4-pentylpiperidine)-1-yl-propyl)-1H-indazole;

1-(3-(4-propylpiperidine)-1-yl-propyl)-1H-indazole;

1-(3-(4-(3-methyl-butyl)-piperidine)-1-yl-propyl)-1H-indazole

1-(3-(4-pentylidene-piperidine)-1-yl-propyl)-1H-indazole;

1-(3-(4-propylidene-piperidine)-1-yl-propyl)-1H-indazole

1-benzo[*b*]thiophen-2-yl-4-(4-butylpiperidin-1-yl)-butan-1-one

4-(4-butylpiperidin-1-yl)-1-(3-methyl-benzofuran-2-yl)-butan-1-one;

4-(4-butylpiperidin-1-yl)-1-(5-fluoro-3-methyl-benzo[*b*]thiophen-2-yl)-butan-1-one;

1-benzofuran-2-yl-4-(4-butylpiperidin-1-yl)-butan-1-one;

1-(3-bromo-benzo[*b*]thiophen-2-yl)-4-(4-butylpiperidin-1-yl)-butan-1-one

1-(3-benzo[*b*]thiophen-2-yl-propyl)-4-butylpiperidine;

1-(3-benzofuran-2-yl-propyl)-4-butylpiperidine;

4-butyl-1-[3-(3-methyl-benzofuran-2-yl)-propyl]-piperidine;

4-butyl-1-[3-(5-fluoro-3-methyl-benzo[*b*]thiophen-2-yl)-propyl]-piperidine;

2-(3-iodo-propyl)-benzo[*b*]thiophene;

1-(3-benzo[*b*]thiophen-2-yl-propyl)-4-methylpiperidine

1-(3-benzo[*b*]thiophen-2-yl-propyl)-4-benzylpiperidine;

1-(3-benzo[*b*]thiophen-2-yl-propyl)-4-(2-methoxy-phenyl)-piperidine;

1-[3-(4-butylpiperidin-1-yl)-propyl]-1H-benzotriazole;

1-[3-(4-butylpiperidin-1-yl)-propyl]-1*H*-indole-3-carbaldehyde;

{1-[3-(4-butylpiperidin-1-yl)-propyl]-1H-indol-3-yl}-methanol;

1-[3-(4-butylpiperidin-1-yl)-propyl]-2-phenyl-1*H*-benzoimidazole;

1-[3-(4-butylpiperidin-1-yl)-propyl]-3-chloro-1*H*-indazole;

1-[3-(4-butylpiperidin-1-yl)-propyl]-6-nitro-1*H*-indazole;

3-[2-(4-butylpiperidin-1-yl)-ethoxy]-benzo[*d*]isoxazol;

3-[3-(4-butyl-piperidin-1-yl)-propyl]- 1*H*-indole hydrochloride;

1H-indazole-3-carboxylic acid (2-(4-butylpiperidin)-1-yl-ethyl)-amide;

1-[3-(4-butylpiperidin-1-yl)-propyl]-5-nitro-1H-indazole;

1-[3-(4-butyl-piperidin-1-yl)-propyl]-2-methyl-1*H*-indole;

1-{1-[3-(4-butyl-piperidin-1-yl)-propyl]-1H-indol-3-yl}-ethanone;

{1-[3-(4-butyl-piperidin-1-yl)-propyl]-1H-indol-3-yl}-acetonitrile;

1-[3-(4-butyl-piperidin-1-yl)-propyl]-1*H*-indole -3-carbonitrile;

1-[3-(4-butyl-piperidin-1-yl)-propyl]-5,6-dimethyl-1*H*-benzoimidazole;

1-[3-(4-butyl-piperidin-1-yl)-propyl]-5(6)-dimethyl-1*H*-benzoimidazole;

1-[3-(4-butyl-piperidin-1-yl)-propyl]-5-methoxy-1*H*-benzoimidazole;

{1-[3-(4-butyl-piperidin-1-yl)-propyl]-1H-benzoimidazol-2-yl}-methanol;

1-[3-(4-butyl-piperidin-1-yl)-propyl]-2-trifuoromethyl-1*H*-benzoimidazole;

3-[3-(4-butyl-piperidine-1-yl)-propyl]-1H-indazole, HCl;

3-[3-(4-butyl-piperidine-1-yl)-propyl]-5-nitro-1H-indazole;

3-[3-(4-butyl-piperidine-1-yl)-propyl]-5,7-dinitro-1H-indazole;

3-[3-(4-butyl-piperidin-1-yl)-propyl]-benzo[*d*]isothiazole;

3-[3-(4-butyl-piperidin-1-yl)-propyl]-5-methoxy-1H-indazole;

3-[3-(4-butyl-piperidin-1-yl)-propyl]-4-methoxy-1H-indazole

3-[3-(4-butyl-piperidin-1-yl)-propyl]-6-methoxy-1H-indazole;

3-[3-(4-butyl-piperidin-1-yl)-propyl]-1H-indazole-4-ol;

3-[3-(4-butyl-piperidin-1-yl)-propyl]-1H-indazole-6-ol; or

3-[3-(4-butyl-piperidin-1-yl)-propyl]-1H-indazole-5-ol.

35-76. (CANCELED)